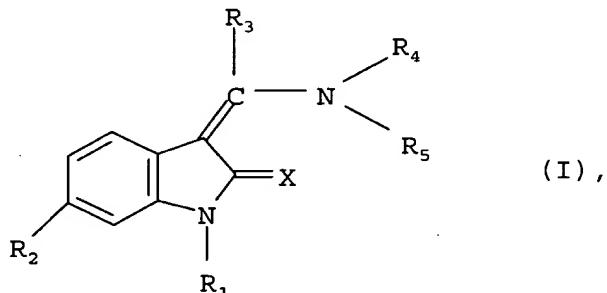


AMENDMENTS TO THE CLAIMS

Claim 1 (Previously amended):

A compound of the formula I



(I),

wherein:

X denotes an oxygen atom,

R₁ denotes a hydrogen atom or a C₁₋₄-alkoxycarbonyl or C₂₋₄-alkanoyl group,

C 1

R₂ denotes a carboxy group, a straight-chain or branched C₁₋₆-alkoxy-carbonyl group, a C₄₋₇-cycloalkoxy-carbonyl or an aryloxycarbonyl group,

a straight-chain or branched C₁₋₆-alkoxy-carbonyl group, which is terminally substituted in the alkyl moiety by a phenyl, heteroaryl, carboxy, C₁₋₃-alkoxy-carbonyl, aminocarbonyl, C₁₋₃-alkylamino-carbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl group,

a straight-chain or branched C₂₋₆-alkoxy-carbonyl group, which is terminally substituted in the alkyl moiety by a chlorine atom or a hydroxy, C₁₋₃-alkoxy, amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

an aminocarbonyl or methylaminocarbonyl group, an ethylaminocarbonyl group optionally substituted in the 2 position of the ethyl group by a hydroxy or C₁₋₃-alkoxy group, |

R₃ denotes a hydrogen atom, a C₁₋₆-alkyl, C₃₋₇-cycloalkyl, trifluoromethyl or heteroaryl group,

a phenyl or naphthyl group, a phenyl or naphthyl group mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom, by a trifluoromethyl, C₁₋₃-alkyl or C₁₋₃-alkoxy group, whilst in the event of disubstitution the substituents may be identical or different and wherein the abovementioned unsubstituted as well as the mono- and disubstituted phenyl and naphthyl groups may additionally be substituted

by a hydroxy, hydroxy-C₁₋₃-alkyl or C₁₋₃-alkoxy-C₁₋₃-alkyl group,

by a cyano, carboxy, carboxy-C₁₋₃-alkyl, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylamino-carbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl group,

by a nitro group,

C 1 by an amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino or amino-C₁₋₃-alkyl group,

by a C₁₋₃-alkylcarbonylamino, N-(C₁₋₃-alkyl)-C₁₋₃-alkyl-carbonylamino, C₁₋₃-alkylcarbonylamino-C₁₋₃-alkyl, N-(C₁₋₃-alkyl)-C₁₋₃-alkylcarbonylamino-C₁₋₃-alkyl, C₁₋₃-alkyl-sulphonylamino, C₁₋₃-alkylsulphonylamino-C₁₋₃-alkyl, N-(C₁₋₃-alkyl)-C₁₋₃-alkylsulphonylamino-C₁₋₃-alkyl or aryl-C₁₋₃-alkylsulphonylamino group,

by a cycloalkylamino, cycloalkyleneimino, cycloalkyleneiminocarbonyl, cycloalkyleneimino-C₁₋₃-alkyl, cycloalkyleneiminocarbonyl-C₁₋₃-alkyl or cycloalkyleneiminosulphonyl-C₁₋₃-alkyl group having 4 to 7 ring members in each case, whilst in each case the methylene group in position 4 of a 6- or 7-membered cycloalkyleneimino group may be replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH or -N(C₁₋₃-alkyl) group,

or by a heteroaryl or heteroaryl-C₁₋₃-alkyl group,

R₄ denotes a C₃₋₇-cycloalkyl group,

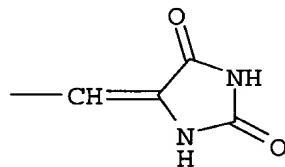
whilst the methylene group in the 4 position of a 6- or 7-membered cycloalkyl group may be substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or replaced by an -NH or -N(C₁₋₃-alkyl) group,

or a phenyl group substituted by the group R₆, which may additionally be mono- or disubstituted by fluorine, chlorine, bromine or iodine atoms, by C₁₋₅-alkyl, trifluoromethyl, hydroxy, C₁₋₃-alkoxy, carboxy, C₁₋₃-alkoxycarbonyl, amino, acetylarnino, C₁₋₃-alkyl-sulphonylarnino, aminocarbonyl, C₁₋₃-alkyl-aminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, aminosulphonyl, C₁₋₃-alkyl-aminosulphonyl, di-(C₁₋₃-alkyl)-aminosulphonyl, nitro or cyano groups, wherein the substituents may be identical or different and wherein

R₆ denotes a hydrogen, fluorine, chlorine, bromine or iodine atom,

a cyano, nitro, amino, C₁₋₅-alkyl, C₃₋₇-cycloalkyl, trifluoromethyl, phenyl, tetrazolyl or heteroaryl group,

the group of formula



wherein the hydrogen atoms bound to a nitrogen atom may in each case be replaced independently of one another by a C₁₋₃-alkyl group,

a C₁₋₃-alkoxy group, a C₁₋₃-alkoxy-C₁₋₃-alkoxy, phenyl-C₁₋₃-alkoxy, amino-C₂₋₃-alkoxy, C₁₋₃-alkylarnino-C₂₋₃-alkoxy, di-(C₁₋₃-alkyl)-arnino-C₂₋₃-alkoxy, phenyl-C₁₋₃-alkylarnino-C₂₋₃-alkoxy, N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylarnino-C₂₋₃-alkoxy, C₅₋₇-cycloalkyleneimino-C₂₋₃-alkoxy or C₁₋₃-alkylmercapto group,

a carboxy, C₁₋₄-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylarnino-carbonyl, N-(C₁₋₅-alkyl)-C₁₋₃-alkylaminocarbonyl, phenyl-C₁₋₃-alkylarnino-carbonyl, N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylarnino-carbonyl, piperazinocarbonyl or N-(C₁₋₃-alkyl)-piperazinocarbonyl group,

a C₁₋₃-alkylaminocarbonyl or N-(C₁₋₅-alkyl)-C₁₋₃-alkylaminocarbonyl group wherein an alkyl moiety is substituted by a carboxy or C₁₋₃-alkoxycarbonyl group or in the 2 or 3 position by a di-(C₁₋₃-alkyl)-amino, piperazino, N-(C₁₋₃-alkyl)-piperazino or a 4- to 7-membered cycloalkyleneimino group,

a C₃₋₇-cycloalkyl-carbonyl group,

wherein the methylene group in the 4 position of the 6- or 7-membered cycloalkyl moiety may be substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or replaced by an -NH or -N(C₁₋₃-alkyl) group,

a 4- to 7-membered cycloalkyleneimino group wherein

C (

a methylene group linked to the imino group may be replaced by a carbonyl or sulphonyl group or

the cycloalkylene moiety may be fused to a phenyl ring or

one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl group and/or

in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be substituted by a carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, phenyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino group or

may be replaced by an oxygen or sulphur atom, by a sulphanyl, sulphonyl, -NH, -N(C₁₋₃-alkyl), -N(phenyl), -N(C₁₋₃-alkyl-carbonyl) or -N(benzoyl) group,

a C₁₋₄-alkyl group substituted by the group R₇, wherein

R₇ denotes a C₃₋₇-cycloalkyl group,

whilst the methylene group in the 4 position of a 6- or 7-membered cycloalkyl group may be substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or replaced by an -NH or -N(C₁₋₃-alkyl) group or

in a 5- to 7-membered cycloalkyl group a -(CH₂)₂ group may be replaced by a -CO-NH group, a -(CH₂)₃ group may be replaced by a -NH-CO-NH or -CO-NH-CO group or a -(CH₂)₄ group may be replaced by a -NH-CO-NH-CO group, whilst in each case a hydrogen atom bound to a nitrogen atom may be replaced by a C₁₋₃-alkyl group,

an aryl or heteroaryl group,

a hydroxy or C₁₋₃-alkoxy group,

(1)

an amino, C₁₋₇-alkylamino, di-(C₁₋₇-alkyl)-amino, phenylamino, N-phenyl-C₁₋₃-alkyl-amino, phenyl-C₁₋₃-alkylamino, N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino or di-(phenyl-C₁₋₃-alkyl)-amino group,

an ω -hydroxy-C₂₋₃-alkyl-amino, N-(C₁₋₃-alkyl)- ω -hydroxy-C₂₋₃-alkyl-amino, di-(ω -hydroxy-C₂₋₃-alkyl)-amino, di-(ω -(C₁₋₃-alkoxy)-C₂₋₃-alkyl)-amino or N-(dioxolan-2-yl)-C₁₋₃-alkyl-amino group,

a C₁₋₃-alkylcarbonylamino-C₂₋₃-alkyl-amino or C₁₋₃-alkylcarbonylamino-C₂₋₃-alkyl-N-(C₁₋₃-alkyl)-amino group,

a C₁₋₃-alkylsulphonylamino, N-(C₁₋₃-alkyl)-C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylsulphonylamino-C₂₋₃-alkyl-amino or C₁₋₃-alkylsulphonylamino-C₂₋₃-alkyl-N-(C₁₋₃-alkyl)-amino group,

a hydroxycarbonyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-hydroxycarbonyl-C₁₋₃-alkyl-amino group,

a guanidino group wherein one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl group,

a group of formula



wherein

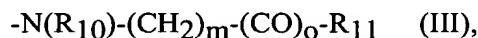
(1)

R₈ denotes a hydrogen atom or a C₁₋₃-alkyl group,

n denotes one of the numbers 0, 1, 2 or 3 and

R₉ denotes an amino, C₁₋₄-alkylamino, di-(C₁₋₄-alkyl)-amino, phenylamino, N-(C₁₋₄-alkyl)-phenylamino, benzylamino, N-(C₁₋₄-alkyl)-benzylamino or C₁₋₄-alkoxy group, a 4- to 7-membered cycloalkyleneimino group, whilst in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH, -N(C₁₋₃-alkyl), -N(phenyl), -N(C₁₋₃-alkyl-carbonyl) or -N(benzoyl) group, or, if n denotes one of the numbers 1, 2 or 3, it may also denote a hydrogen atom,

a group of formula



wherein

R_{10} denotes a hydrogen atom, a C_{1-3} -alkyl group, a C_{1-3} -alkylcarbonyl, arylcarbonyl, phenyl- C_{1-3} -alkyl-carbonyl, C_{1-3} -alkylsulphonyl, arylsulphonyl or phenyl- C_{1-3} -alkylsulphonyl group,

m denotes one of the numbers 1, 2, 3 or 4,

o denotes the number 1 or, if m denotes one of the numbers 2, 3 or 4, o may also denote the number 0 and

R_{11} denotes an amino, C_{1-4} -alkylamino, di-(C_{1-4} -alkyl)-amino, phenylamino, N -(C_{1-4} -alkyl)-phenylamino, benzylamino, N -(C_{1-4} -alkyl)-benzylamino, C_{1-4} -alkoxy or C_{1-3} -alkoxy- C_{1-3} -alkoxy group, a di-(C_{1-4} -alkyl)-amino- C_{1-3} -alkylamino group optionally substituted in the 1 position by a C_{1-3} -alkyl group or a 4- to 7-membered cycloalkyleneimino group, wherein the cycloalkylene moiety may be fused to a phenyl ring or in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH, - $N(C_{1-3}$ -alkyl), -N(phenyl), - $N(C_{1-3}$ -alkyl-carbonyl) or -N(benzoyl) group,

a C_{4-7} -cycloalkylamino, C_{4-7} -cycloalkyl- C_{1-3} -alkylamino or C_{4-7} -cycloalkenylamino group wherein position 1 of the ring is not involved in the double bond and wherein the abovementioned groups may each additionally be substituted at the amino-nitrogen atom by a C_{5-7} -cycloalkyl, C_{2-4} -alkenyl or C_{1-4} -alkyl group,

a 4- to 7-membered cycloalkyleneimino group, wherein

the cycloalkylene moiety may be fused to a phenyl group or to an oxazolo, imidazolo, thiazolo, pyridino, pyrazino or pyrimidino group optionally substituted by a fluorine, chlorine, bromine or iodine atom, by a nitro, C_{1-3} -alkyl, C_{1-3} -alkoxy or amino group, and/or

one or two hydrogen atoms may each be replaced by a C_{1-3} -alkyl, C_{5-7} -cycloalkyl or phenyl group and/or

the methylene group in the 3 position of a 5-membered cycloalkyleneimino group may be substituted by a hydroxy, hydroxy-C₁₋₃-alkyl, C₁₋₃-alkoxy or C₁₋₃-alkoxy-C₁₋₃-alkyl group,

the methylene group in the 3 or 4 position of a 6- or 7-membered cycloalkyleneimino group may in each case be substituted by a hydroxy, hydroxy-C₁₋₃-alkyl, C₁₋₃-alkoxy, C₁₋₃-alkoxy-C₁₋₃-alkyl, carboxy, C₁₋₄-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, phenyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkyl-amino group or

may be replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH, -N(C₁₋₃-alkyl-), -N(phenyl), -N(phenyl-C₁₋₃-alkyl-), -N(C₁₋₃-alkyl-carbonyl-), -N(C₁₋₄-hydroxy-carbonyl-), -N(C₁₋₄-alkoxy-carbonyl-), -N(benzoyl-) or -N(phenyl-C₁₋₃-alkyl-carbonyl-) group,

wherein a methylene group linked to an imino-nitrogen atom of the cycloalkyleneimino group may be replaced by a carbonyl or sulphonyl group or in a 5- to 7-membered monocyclic cycloalkyleneimino group or a cycloalkyleneimino group fused to a phenyl group the two methylene groups linked to the imino-nitrogen atom may each be replaced by a carbonyl group,

or R₆ denotes a C₁₋₄-alkyl group which is substituted by a carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl group or by a 4- to 7-membered cycloalkyleneiminocarbonyl group,

an N-(C₁₋₃-alkyl)-C₂₋₄-alkanoylamino group which is additionally substituted in the alkyl moiety by a carboxy or C₁₋₃-alkoxycarbonyl group,

a group of formula



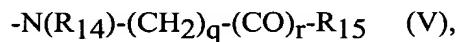
wherein

R_{12} denotes a hydrogen atom, a C_{1-6} -alkyl or C_{3-7} -cycloalkyl group or a C_{1-3} -alkyl group terminally substituted by a phenyl, heteroaryl, trifluoromethyl, hydroxy, C_{1-3} -alkoxy, aminocarbonyl, C_{1-4} -alkylamino-carbonyl, di-(C_{1-4} -alkyl)-amino-carbonyl, C_{1-3} -alkyl-carbonyl, C_{1-3} -alkyl-sulphonylamino, $N-(C_{1-3}$ -alkyl)- C_{1-3} -alkyl-sulphonylamino, C_{1-3} -alkyl-aminosulphonyl or di-(C_{1-3} -alkyl)-aminosulphonyl group and

p denotes one of the numbers 0, 1, 2 or 3 and

C_1 assumes the meanings of the abovementioned group R_7 , or, if p denotes one of the numbers 1, 2 or 3, it may also denote a hydrogen atom,

a group of formula



wherein

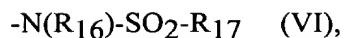
R_{14} denotes a hydrogen atom, a C_{1-4} -alkyl group, a C_{1-3} -alkylcarbonyl, arylcarbonyl, phenyl- C_{1-3} -alkylcarbonyl, heteroarylcarbonyl, heteroaryl- C_{1-3} -alkylcarbonyl, C_{1-4} -alkylsulphonyl, arylsulphonyl, phenyl- C_{1-3} -alkylsulphonyl, heteroarylsulphonyl or heteroaryl- C_{1-3} -alkyl-sulphonyl group,

q denotes one of the numbers 1, 2, 3 or 4,

r denotes the number 1 or, if q is one of the numbers 2, 3 or 4, it may also denote the number 0 and

R_{15} assumes the meanings of the abovementioned group R_7 ,

a group of formula



wherein

R_{16} denotes a hydrogen atom or a C_{1-4} -alkyl group optionally terminally substituted by a cyano, trifluoromethyl-carbonylamino or $\text{N}-(\text{C}_{1-3}\text{-alkyl})$ -trifluoromethyl-carbonyl-amino group and

R_{17} denotes a C_{1-3} -alkyl group,

\cap (an amino group substituted by a di- $(\text{C}_{1-3}\text{-alkyl})$ -amino- C_{1-3} -alkyl-carbonyl or di- $(\text{C}_{1-3}\text{-alkyl})$ -amino- C_{1-3} -alkyl-sulphonyl group and a di- $(\text{C}_{1-3}\text{-alkyl})$ -aminocarbonyl- C_{1-3} -alkyl group,

or an $\text{N}-(\text{C}_{1-3}\text{-alkyl})$ - C_{1-5} -alkylsulphonylamino or $\text{N}-(\text{C}_{1-3}\text{-alkyl})$ -phenylsulphonylamino group wherein the alkyl moiety is additionally substituted by a cyano or carboxy group,

wherein all the single-bonded or fused phenyl groups contained in the groups mentioned under R_6 may be mono- or disubstituted by fluorine, chlorine, bromine or iodine atoms, by C_{1-5} -alkyl, trifluoromethyl, hydroxy, C_{1-3} -alkoxy, carboxy, C_{1-3} -alkoxycarbonyl, aminocarbonyl, C_{1-4} -alkylamino-carbonyl, di- $(\text{C}_{1-4}\text{-alkyl})$ -amino-carbonyl, aminosulphonyl, C_{1-3} -alkyl-aminosulphonyl, di- $(\text{C}_{1-3}\text{-alkyl})$ -aminosulphonyl, C_{1-3} -alkyl-sulphonylamino, nitro or cyano groups, wherein the substituents may be identical or different, or two adjacent hydrogen atoms of the phenyl groups may be replaced by a methylenedioxy group,

and

R_5 denotes a hydrogen atom or a C_{1-3} -alkyl group,

wherein by an aryl group is meant a phenyl or naphthyl group optionally mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom, by a cyano, trifluoromethyl, nitro, carboxy, aminocarbonyl, C₁₋₃-alkyl or C₁₋₃-alkoxy group and

by a heteroaryl group is meant a monocyclic 5- or 6-membered heteroaryl group optionally substituted by a C₁₋₃-alkyl group in the carbon skeleton, wherein

the 6-membered heteroaryl group contains one, two or three nitrogen atoms and

the 5-membered heteroaryl group contains an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group, an oxygen or sulphur atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group or an oxygen or sulphur atom and additionally a nitrogen atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group and two nitrogen atoms,

and moreover a phenyl ring may be fused to the abovementioned monocyclic heterocyclic groups via two adjacent carbon atoms and the bonding takes place via a nitrogen atom or via a carbon atom of the heterocyclic moiety or a fused phenyl ring,

some or all of the hydrogen atoms in the abovementioned alkyl and alkoxy groups or in the alkyl moieties contained in the above-defined groups of formula I may be replaced by fluorine atoms,

and wherein any carboxy group contained in the abovementioned groups may be replaced by a tert.butoxycarbonyl precursor group,

and wherein a hydrogen atom bound to a nitrogen atom may each be replaced by hydroxyl, benzoyl, pyridinoyl, formyl, acetyl, propionyl, butanoyl, pentanoyl, hexanoyl, allyloxycarbonyl, methoxycarbonyl, ethoxycarbonyl, propoxycarbonyl, isopropoxycarbonyl, butoxycarbonyl, tert.butoxycarbonyl, pentoxy carbonyl, hexyloxycarbonyl, octyloxycarbonyl, nonyloxycarbonyl, decyloxycarbonyl, undecyloxycarbonyl, dodecyloxycarbonyl,

hexadecyloxycarbonyl, benzyloxycarbonyl, phenylethoxycarbonyl, phenylpropoxycarbonyl, C₁₋₃-alkylsulphonyl-C₂₋₄-alkoxycarbonyl, C₁₋₃-alkoxy-C₂₋₄-alkoxy-C₂₋₄-alkoxycarbonyl or an R_eCO-O-(R_fCR_g)-O-CO group wherein

R_e denotes a C₁₋₈-alkyl, C₅₋₇-cycloalkyl, phenyl or phenyl-C₁₋₃-alkyl group,

R_f denotes a hydrogen atom, a C₁₋₃-alkyl, C₅₋₇-cycloalkyl or phenyl group and

R_g denotes a hydrogen atom, a C₁₋₃-alkyl or R_eCO-O-(R_fCR_g)-O group wherein R_e to R_g are as hereinbefore defined,

or wherein an amino nitrogen may form part of a phthalimido group,

C 1

or a tautomer or salt thereof.

Claim 2 (Previously amended): A compound of the formula I according to claim 1, wherein:

R₁ and R₃ are as defined in claim 1,

X denotes an oxygen atom,

R₂ denotes a carboxy group, a straight-chain or branched C₁₋₆-alkoxy-carbonyl group, a C₅₋₇-cycloalkoxycarbonyl or a phenoxy carbonyl group,

a straight-chain or branched C₁₋₃-alkoxy-carbonyl group, which is terminally substituted in the alkyl moiety by a phenyl, heteroaryl, carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl group,

a straight-chain or branched C₂₋₃-alkoxy-carbonyl group, which is terminally substituted in the alkyl moiety by a chlorine atom, by a hydroxy, C₁₋₃-alkoxy, amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

an aminocarbonyl or methylaminocarbonyl group, an ethylaminocarbonyl group optionally substituted in the 2 position of the ethyl group by a hydroxy or C₁₋₃-alkoxy group, |

R₄ denotes a C₃₋₇-cycloalkyl group,

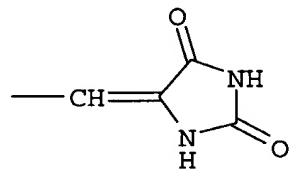
whilst the methylene group in the 4 position of a 6- or 7-membered cycloalkyl group may be substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or replaced by an -NH or -N(C₁₋₃-alkyl) group,

or a phenyl group substituted by the group R₆, which may additionally be mono- or disubstituted by fluorine, chlorine or bromine atoms, by C₁₋₃-alkyl, trifluoromethyl, hydroxy, C₁₋₃-alkoxy, carboxy, C₁₋₃-alkoxycarbonyl, amino, acetylarnino, aminocarbonyl, C₁₋₃-alkyl-aminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, nitro or cyano groups, wherein the substituents may be identical or different and wherein

R₆ denotes a hydrogen, fluorine, chlorine, bromine or iodine atom,

a cyano, nitro, amino, C₁₋₅-alkyl, C₃₋₇-cycloalkyl, trifluoromethyl, phenyl, tetrazolyl or heteroaryl group,

the group of formula



wherein a hydrogen atom bound to the nitrogen atom may be replaced by a C₁₋₃-alkyl group,

a C₁₋₃-alkoxy group, an amino-C₂₋₃-alkoxy, C₁₋₃-alkylamino-C₂₋₃-alkoxy, di-(C₁₋₃-alkyl)-amino-C₂₋₃-alkoxy, phenyl-C₁₋₃-alkylamino-C₂₋₃-alkoxy, N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino-C₂₋₃-alkoxy, pyrrolidino-C₂₋₃-alkoxy, piperidino-C₂₋₃-alkoxy or C₁₋₃-alkylmercapto group,

a carboxy, C₁₋₄-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylamino-carbonyl, phenyl-C₁₋₃-alkylamino-carbonyl or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino-carbonyl group,

a C₃₋₇-cycloalkyl-carbonyl group,

wherein the methylene group in the 4 position of the 6- or 7-membered cycloalkyl moiety may be replaced by an -NH or -N(C₁₋₃-alkyl) group,

a 4- to 7-membered cycloalkyleneimino group, wherein

a methylene group linked to the imino group may be replaced by a carbonyl or sulphonyl group or

one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl group and/or

C 1

in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be substituted by a carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, phenyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino group or

may be replaced by an oxygen or sulphur atom, by a sulphanyl, sulphonyl, -NH or -N(C₁₋₃-alkyl) group,

a C₁₋₄-alkyl group terminally substituted by the group R₇, wherein

R₇ denotes a C₅₋₇-cycloalkyl group,

whilst the methylene group in the 4 position of a 6- or 7-membered cycloalkyl group may be replaced by an -NH or -N(C₁₋₃-alkyl) group or

in a 5- to 7-membered cycloalkyl group a -(CH₂)₂ group may be replaced by a -CO-NH group, a -(CH₂)₃ group may be replaced by a -NH-CO-NH- or a -(CH₂)₄

group may be replaced by a -NH-CO-NH-CO group, whilst in each case a hydrogen atom bound to a nitrogen atom may be replaced by a C₁₋₃-alkyl group,

a phenyl or heteroaryl group,

a hydroxy or C₁₋₃-alkoxy group,

an amino, C₁₋₆-alkylamino, di-(C₁₋₆-alkyl)-amino, phenylamino, N-phenyl-C₁₋₃-alkyl-amino, phenyl-C₁₋₃-alkylamino, N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino or di-(phenyl-C₁₋₃-alkyl)-amino group,

a ω -hydroxy-C₂₋₃-alkyl-amino, N-(C₁₋₃-alkyl)- ω -hydroxy--C₂₋₃-alkyl-amino, di-(ω -hydroxy-C₂₋₃-alkyl)-amino, di-(ω -(C₁₋₃-alkoxy)-C₂₋₃-alkyl)-amino or N-(dioxolan-2-yl)-C₁₋₃-alkyl-amino group,

a C₁₋₃-alkylcarbonylamino-C₂₋₃-alkyl-amino or C₁₋₃-alkylcarbonylamino-C₂₋₃-alkyl-N-(C₁₋₃-alkyl)-amino group,

a C₁₋₃-alkylsulphonylamino, N-(C₁₋₃-alkyl)-C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylsulphonylamino--C₂₋₃-alkyl-amino or C₁₋₃-alkylsulphonylamino-C₂₋₃-alkyl-N-(C₁₋₃-alkyl)-amino group,

a hydroxycarbonyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-hydroxycarbonyl-C₁₋₃-alkyl-amino group

a guanidino group wherein a hydrogen atom may be replaced by a C₁₋₃-alkyl group,

a group of formula

$-N(R_8)-CO-(CH_2)_n-R_9$ (II),

wherein

R_8 denotes a hydrogen atom or a C_{1-3} -alkyl group,

n denotes one of the numbers 0, 1, 2 or 3 and

R_9 denotes an amino, C_{1-3} -alkylamino, di- $(C_{1-3}$ -alkyl)-amino, phenylamino, benzylamino or C_{1-4} -alkoxy group, a 5- to 7-membered cycloalkyleneimino group, wherein the methylene group in position 4 of the piperidino group may be replaced by an oxygen or sulphur atom, by an -NH, $-N(C_{1-3}$ -alkyl), $-N(phenyl)$, $-N(C_{1-3}$ -alkyl-carbonyl) or $-N(benzoyl)$ group, or, if n denotes one of the numbers 1, 2 or 3, it may also denote a hydrogen atom,

C_l

a group of formula

$-N(R_{10})-(CH_2)_m-(CO)_o-R_{11}$ (III),

wherein

R_{10} denotes a hydrogen atom, a C_{1-3} -alkyl group, a C_{1-3} -alkylcarbonyl or C_{1-3} -alkylsulphonyl group,

m denotes one of the numbers 1, 2 or 3,

o denotes the number 1 or, if m is one of the numbers 2 or 3, o may also denote the number 0 and

R_{11} denotes an amino, C_{1-3} -alkylamino, di- $(C_{1-3}$ -alkyl)-amino, C_{1-4} -alkoxy or C_{1-3} -alkoxy- C_{1-3} -alkoxy group or a 5- to 7-membered cycloalkyleneimino group, wherein the methylene group in position 4 of the piperidino group may be replaced

by an oxygen or sulphur atom, by an -NH, -N(C₁₋₃-alkyl), -N(phenyl), -N(C₁₋₃-alkyl-carbonyl) or -N(benzoyl) group,

a C₄₋₇-cycloalkylamino or C₄₋₇-cycloalkenylamino group wherein position 1 of the ring is not involved in the double bond,

a 4- to 7-membered cycloalkyleneimino group, wherein

the cycloalkylene moiety may be fused to a phenyl group or

one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl group and/or

the methylene group in position 3 of the pyrrolidino group may be substituted by a hydroxy or C₁₋₃-alkoxy group,

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in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be substituted by a hydroxy, hydroxy-C₁₋₃-alkyl, C₁₋₃-alkoxy, carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylamino-carbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, phenyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino group or

may be replaced by an oxygen or sulphur atom, by a sulphanyl, sulphonyl, -NH, -N(C₁₋₃-alkyl), -N(phenyl), -N(phenyl-C₁₋₃-alkyl), -N(C₁₋₃-alkyl-carbonyl), -N(C₁₋₄-alkoxy-carbonyl), -N(benzoyl) or -N(phenyl-C₁₋₃-alkyl-carbonyl) group,

wherein a methylene group linked to an imino-nitrogen atom of the cycloalkyleneimino group may be replaced by a carbonyl or sulphonyl group or in a 5- to 6-membered monocyclic cycloalkyleneimino group or a cycloalkyleneimino group fused to a phenyl group the two methylene groups linked to the imino-nitrogen atom may each be replaced by a carbonyl group,

or R₆ denotes a C₁₋₄-alkyl group which is terminally substituted by a carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl group or by a 4- to 7-membered cycloalkyleneiminocarbonyl group,

a group of formula



wherein

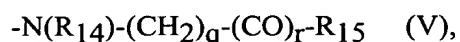
R₁₂ denotes a hydrogen atom, a C₁₋₃-alkyl, C₅₋₇-cycloalkyl, phenyl-C₁₋₃-alkyl or heteroaryl-C₁₋₃-alkyl group and

p denotes one of the numbers 0, 1, 2 or 3 and

C 1

R₁₃ assumes the meanings of the abovementioned group R₇, or, if p denotes one of the numbers 1, 2 or 3, it may also denote a hydrogen atom,

a group of formula



wherein

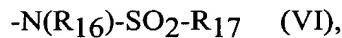
R₁₄ denotes a hydrogen atom, a C₁₋₄-alkyl group, a C₁₋₃-alkylcarbonyl, phenylcarbonyl, phenyl-C₁₋₃-alkylcarbonyl, heteroarylcarbonyl, heteroaryl-C₁₋₃-alkylcarbonyl, C₁₋₄-alkylsulphonyl, phenylsulphonyl, phenyl-C₁₋₃-alkylsulphonyl- heteroarylsulphonyl or heteroaryl-C₁₋₃-alkyl-sulphonyl group,

q denotes one of the numbers 1, 2, 3 or 4,

r denotes the number 1 or, if q is one of the numbers 2, 3 or 4, it may also denote the number 0 and

R₁₅ assumes the meanings of the abovementioned group R₇,

a group of formula



wherein

R₁₆ denotes a hydrogen atom or a C₁₋₄-alkyl group optionally terminally substituted by a cyano, trifluoromethyl-carbonylamino or N-(C₁₋₃-alkyl)-trifluoromethyl-carbonyl-amino group and

R₁₇ denotes a C₁₋₃-alkyl group,

C 1

an amino group substituted by a di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl or di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-sulphonyl group and a di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl group,

wherein all the single-bonded or fused phenyl groups contained in the groups mentioned under R₆ may be mono- or disubstituted by fluorine, chlorine or bromine atoms, by C₁₋₃-alkyl, trifluoromethyl, hydroxy, C₁₋₃-alkoxy, carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkyl-aminocarbonyl, aminosulphonyl, C₁₋₃-alkyl-aminosulphonyl, nitro or cyano groups, wherein the substituents may be identical or different, or two adjacent hydrogen atoms of the phenyl groups may be replaced by a methylenedioxy group, and

R₅ denotes a hydrogen atom or a C₁₋₃-alkyl group,

whilst by a heteroaryl group as mentioned above is meant a pyridinyl, pyrazinyl, pyrimidinyl, pyridazinyl, pyrrolyl, furyl, thienyl, oxazolyl, thiazolyl, pyrazolyl, imidazolyl or triazolyl group optionally substituted in the carbon skeleton by a C₁₋₃-alkyl group wherein a hydrogen atom bound to a nitrogen atom may be replaced by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group

and wherein the 5-membered heteroaryl groups containing at least one imino group are bound via a carbon or nitrogen atom,

a hydrogen atom bound to a nitrogen atom in the abovementioned groups may be replaced by hydroxyl, benzoyl, pyridinoyl, formyl, acetyl, propionyl, butanoyl, pentanoyl, hexanoyl, allyloxycarbonyl, methoxycarbonyl, ethoxycarbonyl, propoxycarbonyl, isopropoxycarbonyl, butoxycarbonyl, tert.butoxycarbonyl, pentoxy carbonyl, hexyloxycarbonyl, octyloxycarbonyl, nonyloxycarbonyl, decyloxycarbonyl, undecyloxycarbonyl, dodecyloxycarbonyl, hexadecyloxycarbonyl, benzyloxycarbonyl, phenylethoxycarbonyl, phenylpropoxycarbonyl, C₁₋₃-alkylsulphonyl-C₂₋₄-alkoxycarbonyl, C₁₋₃-alkoxy-C₂₋₄-alkoxy-C₂₋₄-alkoxycarbonyl or an R_eCO-O-(R_fCR_g)-O-CO group wherein

R_e denotes a C₁₋₈-alkyl, C₅₋₇-cycloalkyl, phenyl or phenyl- C₁₋₃-alkyl group,

R_f denotes a hydrogen atom, a C₁₋₃-alkyl, C₅₋₇-cycloalkyl or phenyl group and

R_g denotes a hydrogen atom, a C₁₋₃-alkyl or R_eCO-O-(R_fCR_g)-O group wherein R_e to R_g are as hereinbefore defined,

or wherein an amino nitrogen may form part of a phthalimido group,

,

and wherein any carboxy group contained in the abovementioned groups may be replaced by a tert.butoxycarbonyl precursor group,

and wherein some or all of the hydrogen atoms in the abovementioned alkyl and alkoxy groups or in the alkyl moieties contained in the above-defined groups of formula I may be replaced by fluorine atoms,

or a tautomer or salt thereof.

Claim 3 (Original): A compound of the formula I according to claim 1, wherein:

X denotes an oxygen atom,

R₁ denotes a hydrogen atom,

R₂ denotes a carboxy group, a straight-chain or branched C₁₋₄-alkoxycarbonyl group or a phenoxy carbonyl group,

a straight-chain or branched C₁₋₃-alkoxy-carbonyl group, which is terminally substituted in the alkyl moiety by a phenyl, carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl group,

a straight-chain or branched C₂₋₃-alkoxy-carbonyl group which is terminally substituted in the alkyl moiety by a hydroxy, C₁₋₃-alkoxy, amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

C l
an aminocarbonyl or methylaminocarbonyl group, an ethylaminocarbonyl group optionally substituted in the 2 position of the ethyl group by a hydroxy or C₁₋₃-alkoxy group or, if R₄ does not denote an aminosulphonyl-phenyl or N-(C₁₋₅-alkyl)-C₁₋₃-alkylaminocarbonyl-phenyl group, it may also denote a di-(C₁₋₂-alkyl)-aminocarbonyl group, |

R₃ denotes a C₁₋₄-alkyl group or a phenyl group which may be substituted by a fluorine, chlorine or bromine atom, by a trifluoromethyl, C₁₋₃-alkyl, hydroxy or C₁₋₃-alkoxy group,

R₄ denotes a C₅₋₆-cycloalkyl group,

wherein the methylene group in position 4 of the cyclohexyl group may be substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or replaced by an -NH or -N(C₁₋₃-alkyl) group,

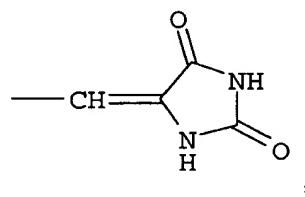
a phenyl group, a phenyl group disubstituted by C₁₋₃-alkyl, C₁₋₃-alkoxy or nitro groups, wherein the substituents may be identical or different, or

a phenyl group substituted by the group R₆, which may additionally be substituted by a fluorine, chlorine or bromine atom or by an amino or nitro group, wherein R₆ denotes a fluorine, chlorine or bromine atom,

a C₁₋₃-alkyl, C₁₋₃-alkoxy, nitro, amino or C₅₋₆-cycloalkyl group,

a pyrrolyl, pyrazolyl, imidazolyl, triazolyl or tetrazolyl group bound via a carbon atom, wherein the abovementioned heteroaromatic groups in the carbon skeleton may be substituted by a C₁₋₃-alkyl group or a hydrogen atom bound to a nitrogen atom may be replaced by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group,

the group of formula



C1

a carboxy, C₁₋₄-alkoxycarbonyl, phenyl-C₁₋₃-alkylamino-carbonyl or C₅₋₇-cycloalkyl-carbonyl group,

a 5 or 6-membered cycloalkyleneimino group, wherein

the methylene group in position 4 of the piperidino group may be replaced by an oxygen or sulphur atom, by an -NH or -N(C₁₋₃-alkyl) group,

an unbranched C₁₋₃-alkyl group terminally substituted by the group R₇, wherein

R₇ denotes a C₅₋₇-cycloalkyl group,

wherein in a 5 or 6-membered cycloalkyl group a -(CH₂)₂ group may be replaced by a -CO-NH group, a -(CH₂)₃ group may be replaced by an -NH-CO-NH- or a

$-(CH_2)_4$ group may be replaced by an -NH-CO-NH-CO group, whilst in each case a hydrogen atom bound to a nitrogen atom may be replaced by a C₁₋₃-alkyl group,

a phenyl or pyridinyl group or a pyrrolyl, pyrazolyl, imidazolyl or triazolyl group bound via a carbon or nitrogen atom, wherein the abovementioned heteroaromatic groups in the carbon skeleton may be substituted by a C₁₋₃-alkyl group or a hydrogen atom bound to a nitrogen atom may be replaced by a C₁₋₃-alkyl group,

a hydroxy or C₁₋₃-alkoxy group,

an amino, C₁₋₆-alkylamino, di-(C₁₋₆-alkyl)-amino, phenylamino, N-phenyl-C₁₋₃-alkylamino, phenyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino group,

1
a ω -hydroxy-C₂₋₃-alkyl-amino, N-(C₁₋₃-alkyl)- ω -hydroxy-C₂₋₃-alkylamino, di-(ω -hydroxy-C₂₋₃-alkyl)-amino or di-(ω -(C₁₋₃-alkoxy)-C₂₋₃-alkyl)-amino group,

a C₁₋₃-alkylcarbonylamino-C₂₋₃-alkyl-amino or C₁₋₃-alkylcarbonylamino-C₂₋₃-alkyl-N-(C₁₋₃-alkyl)-amino group,

a C₁₋₃-alkylsulphonylamino, N-(C₁₋₃-alkyl)-C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylsulphonylamino--C₂₋₃-alkylamino or C₁₋₃-alkylsulphonylamino-C₂₋₃-alkyl-N-(C₁₋₃-alkyl)-amino group,

a hydroxycarbonyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-hydroxycarbonyl-C₁₋₃-alkyl-amino group,

a guanidino group wherein a hydrogen atom may be replaced by a C₁₋₃-alkyl group,

a group of formula



wherein

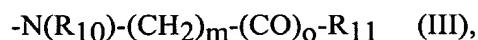
R_8 denotes a hydrogen atom or a C_{1-3} -alkyl group,

n denotes one of the numbers 0, 1, 2 or 3 and

R_9 denotes an amino, C_{1-3} -alkylamino, di-(C_{1-3} -alkyl)-amino or C_{1-4} -alkoxy group, a 5- or 6-membered cycloalkyleneimino group, wherein the methylene group in position 4 of the piperidino group may be replaced by an -NH, -N(C_{1-3} -alkyl) or -N(C_{1-3} -alkyl-carbonyl) group, or, if n denotes one of the numbers 1, 2 or 3, R_9 may also denote a hydrogen atom,

C 1

a group of formula



wherein

R_{10} denotes a hydrogen atom or a C_{1-3} -alkyl group,

m denotes one of the numbers 1, 2 or 3,

o denotes the number 1 or, if m is one of the numbers 2 or 3, o may also denote the number 0 and

R_{11} denotes an amino, C_{1-3} -alkylamino, di-(C_{1-3} -alkyl)-amino, C_{1-4} -alkoxy or methoxy- C_{1-3} -alkoxy group or a 5- or 6-membered cycloalkyleneimino group, wherein the methylene group in position 4 of the piperidino group may be replaced by an -NH, -N(C_{1-3} -alkyl) or -N(C_{1-3} -alkyl-carbonyl) group,

an azetidino, pyrrolidino, piperidino, 2,6-dimethyl-piperidino, 3,5-dimethyl-piperidino or azepino group, wherein

the methylene group in position 3 of the pyrrolidino group may be substituted by a hydroxy group,

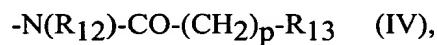
the methylene group in position 4 of the piperidino group may be substituted by a hydroxy, hydroxy-C₁₋₃-alkyl or C₁₋₃-alkoxy group or

may be replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH, -N(C₁₋₃-alkyl), -N(C₁₋₃-alkyl-carbonyl), -N(benzoyl) or -N(phenyl-C₁₋₃-alkyl-carbonyl) group,

wherein a methylene group linked to an imino-nitrogen atom of the pyrrolidino, piperidino or piperazino group may be replaced by a carbonyl group,

or R₆ denotes a straight-chain C₁₋₃-alkyl group which is terminally substituted by a carboxy or C₁₋₃-alkoxy-carbonyl group,

a group of formula



wherein

R₁₂ denotes a hydrogen atom, a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group,

p denotes one of the numbers 0, 1 or 2 and

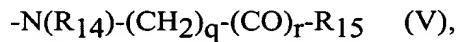
R₁₃ denotes an amino, C₁₋₄-alkylamino, di-(C₁₋₄-alkyl)-amino, benzylamino, N-(C₁₋₃-alkyl)-benzylamino, C₁₋₃-alkoxy-C₁₋₃-alkylamino, N-(C₁₋₃-alkyl)-

C_{1-3} -alkoxy- C_{1-3} -alkylamino, di-(2-methoxy-ethyl)-amino, di-(ω -hydroxy- C_{2-3} -alkyl)-amino or aminocarbonyl-methyl-N-(methyl)-amino group,

a pyrrolyl, pyrazolyl or imidazolyl group bound via a nitrogen atom and optionally substituted by a C_{1-3} -alkyl group,

a pyrrolidino, piperidino, morpholino, thiomorpholino or a piperazino group optionally substituted in the 4 position by a C_{1-3} -alkyl, phenyl- C_{1-3} -alkyl, C_{1-3} -alkylcarbonyl or C_{1-4} -alkoxycarbonyl group or, if n denotes the number 1 or 2, it may also denote a hydrogen atom,

a group of formula



wherein

C1

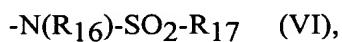
R_{14} denotes a hydrogen atom, a C_{1-4} -alkyl, C_{1-3} -alkyl-carbonyl, phenylcarbonyl, phenyl- C_{1-3} -alkylcarbonyl, furyl-carbonyl, pyridinyl-carbonyl, furyl- C_{1-3} -alkylcarbonyl, pyridinyl- C_{1-3} -alkylcarbonyl, C_{1-4} -alkylsulphonyl, phenylsulphonyl or phenyl- C_{1-3} -alkylsulphonyl group,

q denotes one of the numbers 1, 2 or 3,

r denotes the number 1 or, if q is one of the numbers 2 or 3, it may also denote the number 0 and

R_{15} denotes an amino, C_{1-4} -alkylamino, di-(C_{1-4} -alkyl)-amino, phenylamino, N-(C_{1-4} -alkyl)-phenylamino, benzylamino or N-(C_{1-4} -alkyl)-benzylamino group,

or a group of formula



wherein

R_{16} denotes a hydrogen atom or a C_{1-3} -alkyl group optionally terminally substituted by
a cyano, trifluoromethyl-carbonylamino or
 $N-(C_{1-3}\text{-alkyl})\text{-trifluoromethyl-carbonyl-amino}$ group and

R_{17} denotes a C_{1-3} -alkyl group,

wherein all the single-bonded or fused phenyl groups contained in the groups
mentioned under R_6 may be substituted by a fluorine, chlorine or bromine atom, by
a methyl, trifluoromethyl, methoxy, nitro or cyano group and

R_5 denotes a hydrogen atom,

wherein a hydrogen atom bound to a nitrogen atom in the abovementioned groups may be
replaced by an acetyl or tert.butoxycarbonyl group,

the carboxy groups contained in the abovementioned groups may also be present in the form
of the tert.butoxycarbonyl precursor group,

or a tautomer or salt thereof.

Claim 4 (Original): A compound of the formula I according to claim 1, wherein:

X denotes an oxygen atom,

R_1 and R_5 each denote a hydrogen atom,

R_2 denotes a methoxycarbonyl, ethoxycarbonyl or aminocarbonyl group,

R_3 denotes a phenyl group and

R_4 denotes a phenyl group monosubstituted by the group R_6 , wherein

R_6 denotes an N-methyl-imidazol-2-yl group,

an unbranched C_{1-3} -alkyl group which is terminally substituted by a C_{1-4} -alkylamino, di-(C_{1-4} -alkyl)-amino, piperidino or 2,6-dimethyl-piperidino group,

a group of formula



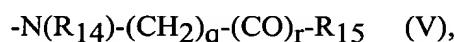
wherein

R_{12} denotes a C_{1-3} -alkyl group,

p denotes one of the numbers 1 or 2 and

R_{13} denotes a di-(C_{1-3} -alkyl)-amino group,

or a group of formula



wherein

R_{14} denotes a C_{1-3} -alkyl-carbonyl or C_{1-3} -alkylsulphonyl group,

q denotes one of the numbers 1, 2 or 3,

r denotes the number 1 or, if q is one of the numbers 2 or 3, r may also denote the number 0 and

R_{15} denotes a di-(C_{1-3} -alkyl)-amino group,

or a tautomer or salt thereof.

Claim 5 (Currently amended): A compound selected from the group consisting of:

- (a) 3-Z-[1-(4-(piperidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-ethoxycarbonyl-2-indolinone,
C 1
- (b) 3-Z-[(1-(4-(piperidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-carbamoyl-2-indolinone,
C 1
- (c) 3-Z-[1-(4-(piperidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
C 1
- (d) 3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-phenyl-methylene]-6-ethoxycarbonyl-2-indolinone,
C 1
- (e) 3-Z-[1-(4-((2,6-dimethyl-piperidin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-ethoxycarbonyl-2-indolinone,
C 1
- (f) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-ethoxycarbonyl-2-indolinone,
C 1
- (g) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-ethoxycarbonyl-2-indolinone,
C 1
- (h) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-ethoxycarbonyl-2-indolinone,
C 1
- (i) 3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
C 1
- (j) 3-Z-[1-(4-(N-acetyl-N-dimethylaminocarbonylmethyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
C 1

- (k) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
C 1
- (l) 3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
and,
- (m) 3-Z-[1-(4-(N-dimethylaminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
and,
- (n) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
and,
- (o) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
and,
- (p) 3-Z-[1-(4-(N-dimethylaminocarbonylmethyl-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
and,
- (q) 3-Z-[1-(4-(N-((2-dimethylamino-ethyl)-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
and,
- (r) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone, and,
- (s) 3-Z-[1-(4-methylaminomethyl-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone, and
- (t) 3-Z-[1-(4-(N-((4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-methoxycarbonyl-2-indolinone,
and,

or a tautomer or salt thereof.

Claim 6 (Original): A physiologically acceptable salt of a compound according to claim 1, 2, 3, 4 or 5.

(l) Claim 7 (Original): A pharmaceutical composition containing a compound according to claim 1, 2, 3 or 4, or a physiologically acceptable salt thereof in accordance with claim 5, together with a pharmaceutically acceptable carrier.

Claim 8 (withdrawn)